



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2017 – 08:28 AM EDT

PDB ID : 4WD7  
Title : Crystal structure of a bacterial Bestrophin homolog from *Klebsiella pneumoniae* by Zn-SAD phasing  
Authors : Yang, T.; Liu, Q.; Hendrickson, W.A.; New York Consortium on Membrane Protein Structure (NYCOMPS)  
Deposited on : unknown  
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029824

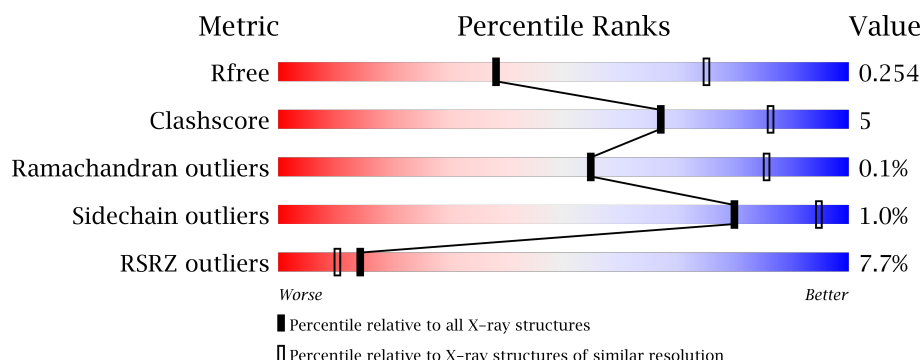
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	301	<div> <div>6%</div> <div>78% 11% 11%</div> </div>
1	B	301	<div> <div>9%</div> <div>80% 10% 10%</div> </div>
1	C	301	<div> <div>6%</div> <div>82% 7% 11%</div> </div>
1	D	301	<div> <div>7%</div> <div>78% 11% 11%</div> </div>
1	E	301	<div> <div>6%</div> <div>76% 12% 10%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10651 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Bestrophin domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2120	1373	359	379	9			
1	B	270	Total	C	N	O	S	0	0	0
			2148	1391	364	384	9			
1	C	269	Total	C	N	O	S	0	0	0
			2110	1362	362	377	9			
1	D	268	Total	C	N	O	S	0	0	0
			2114	1367	359	379	9			
1	E	270	Total	C	N	O	S	0	0	0
			2144	1389	364	382	9			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP S7AS11
A	-1	ASN	-	expression tag	UNP S7AS11
A	0	ALA	-	expression tag	UNP S7AS11
B	-2	SER	-	expression tag	UNP S7AS11
B	-1	ASN	-	expression tag	UNP S7AS11
B	0	ALA	-	expression tag	UNP S7AS11
C	-2	SER	-	expression tag	UNP S7AS11
C	-1	ASN	-	expression tag	UNP S7AS11
C	0	ALA	-	expression tag	UNP S7AS11
D	-2	SER	-	expression tag	UNP S7AS11
D	-1	ASN	-	expression tag	UNP S7AS11
D	0	ALA	-	expression tag	UNP S7AS11
E	-2	SER	-	expression tag	UNP S7AS11
E	-1	ASN	-	expression tag	UNP S7AS11
E	0	ALA	-	expression tag	UNP S7AS11

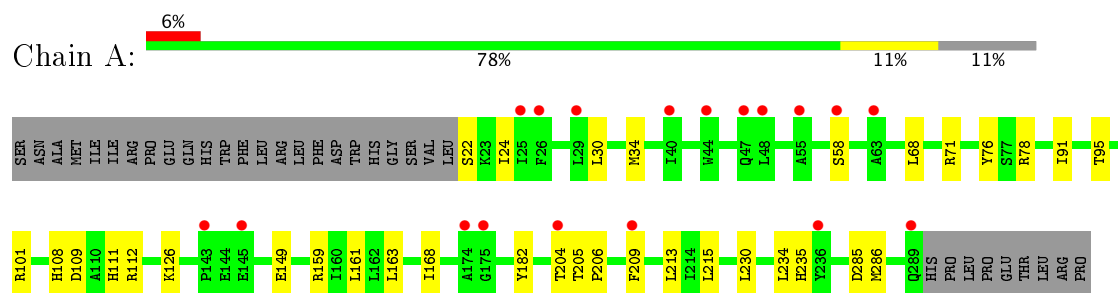
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total 3	Zn 3	0	0
2	A	4	Total 4	Zn 4	0	0
2	D	3	Total 3	Zn 3	0	0
2	C	3	Total 3	Zn 3	0	0
2	E	2	Total 2	Zn 2	0	0

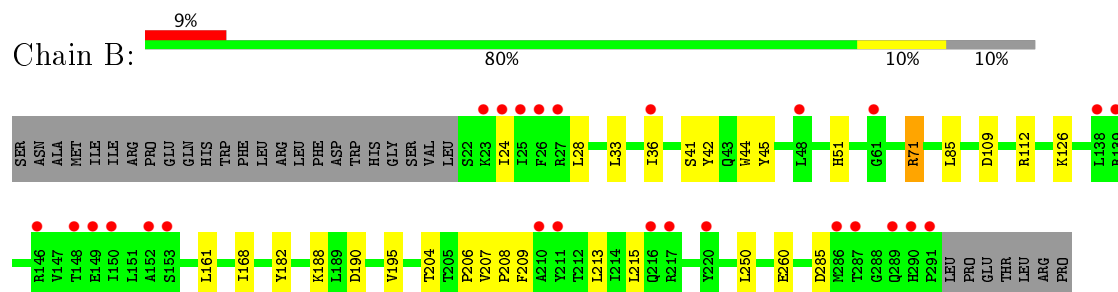
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

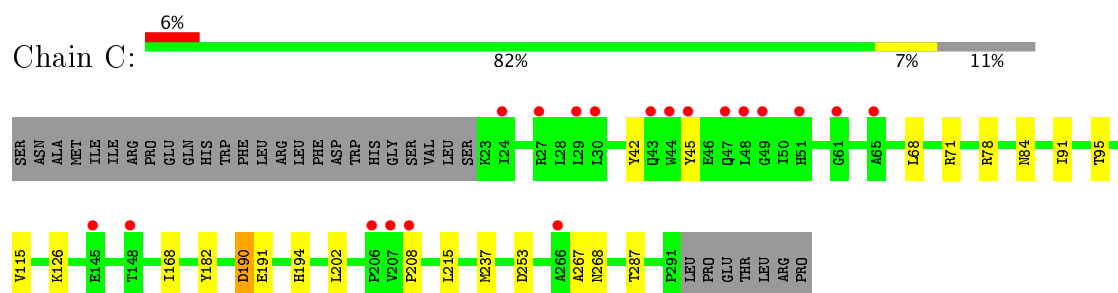
- Molecule 1: Bestrophin domain protein



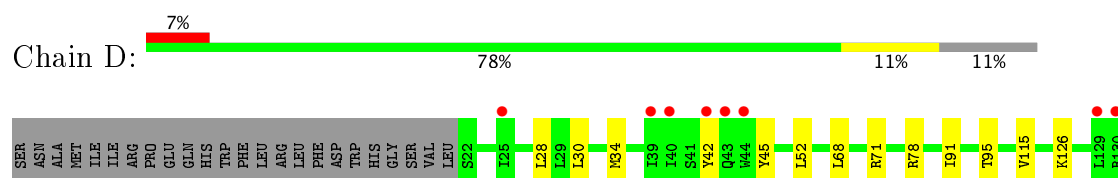
- Molecule 1: Bestrophin domain protein

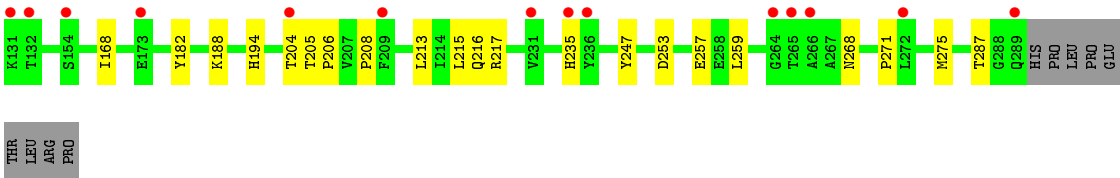


- Molecule 1: Bestrophin domain protein

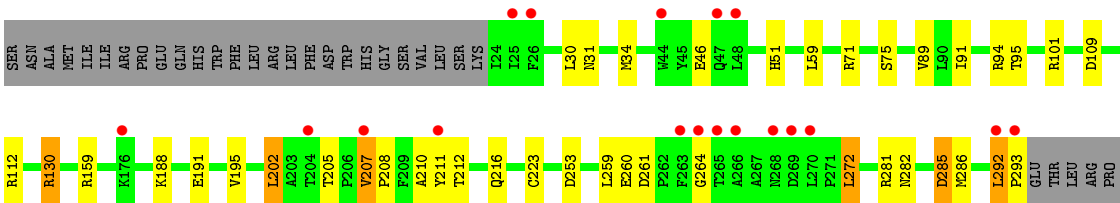
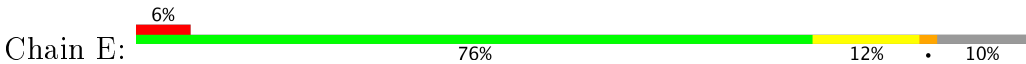


- Molecule 1: Bestrophin domain protein





● Molecule 1: Bestrophin domain protein



GLOBAL-STATISTICS INFOmissingINFO

## 4 Model quality

### 4.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.27	0/2166	0.45	0/2950
1	B	0.26	0/2196	0.44	0/2991
1	C	0.28	0/2155	0.45	0/2937
1	D	0.26	0/2159	0.44	0/2941
1	E	0.29	0/2192	0.54	2/2987 (0.1%)
All	All	0.27	0/10868	0.46	2/14806 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	191	GLU	OE1-CD-OE2	-6.75	115.19	123.30
1	E	130	ARG	NE-CZ-NH2	-5.63	117.48	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	205	THR	Peptide

## 4.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2120	0	2151	24	0
1	B	2148	0	2184	23	0
1	C	2110	0	2134	15	0
1	D	2114	0	2144	25	0
1	E	2144	0	2183	29	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	E	2	0	0	0	0
All	All	10651	0	10796	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (97) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:ASP:OD2	1:E:188:LYS:NZ	2.01	0.94
1:B:208:PRO:HB3	1:D:268:ASN:HD21	1.41	0.83
1:E:109:ASP:OD1	1:E:112:ARG:NH2	2.19	0.75
1:B:190:ASP:OD2	1:D:188:LYS:NZ	2.20	0.74
1:A:68:LEU:HD23	1:A:215:LEU:HD13	1.73	0.71
1:B:109:ASP:OD1	1:B:112:ARG:NH2	2.29	0.65
1:E:207:VAL:HG23	1:E:208:PRO:HD3	1.79	0.63
1:C:71:ARG:NH2	1:C:253:ASP:OD1	2.31	0.62
1:B:24:ILE:HD12	1:B:250:LEU:HB3	1.82	0.62
1:C:68:LEU:HD23	1:C:215:LEU:HD13	1.82	0.61
1:C:78:ARG:NH2	1:C:202:LEU:O	2.31	0.60
1:E:292:LEU:HD22	1:E:293:PRO:HD2	1.82	0.60
1:D:216:GLN:NE2	1:D:257:GLU:OE2	2.34	0.60
1:D:71:ARG:NH2	1:D:253:ASP:OD1	2.32	0.59
1:A:22:SER:HB3	1:A:24:ILE:HG22	1.85	0.58
1:E:31:ASN:ND2	1:E:223:CYS:O	2.36	0.57
1:B:71:ARG:NH1	1:B:260:GLU:OE2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:94:ARG:HH21	1:E:282:ASN:HD21	1.53	0.56
1:C:208:PRO:HG2	1:E:259:LEU:HD21	1.89	0.55
1:C:115:VAL:HG21	1:C:287:THR:HG21	1.88	0.54
1:A:209:PHE:O	1:A:213:LEU:HD13	2.08	0.53
1:B:208:PRO:HG2	1:D:259:LEU:HD21	1.91	0.53
1:A:76:TYR:CE1	1:E:207:VAL:HG21	2.44	0.53
1:A:159:ARG:NH2	1:B:285:ASP:OD2	2.41	0.52
1:D:78:ARG:NH2	1:D:205:THR:O	2.43	0.52
1:A:78:ARG:NH1	1:A:205:THR:O	2.43	0.52
1:A:209:PHE:CZ	1:A:213:LEU:HD11	2.45	0.52
1:C:190:ASP:O	1:C:194:HIS:HD2	1.92	0.52
1:A:109:ASP:OD1	1:A:112:ARG:NH2	2.43	0.52
1:D:115:VAL:HG21	1:D:287:THR:HG21	1.93	0.51
1:A:235:HIS:CD2	1:E:51:HIS:HD2	2.29	0.50
1:E:130:ARG:HG3	1:E:130:ARG:HH11	1.77	0.50
1:D:68:LEU:HD23	1:D:215:LEU:HD13	1.94	0.49
1:B:207:VAL:HG21	1:B:260:GLU:CD	2.31	0.49
1:E:281:ARG:O	1:E:285:ASP:HB2	2.12	0.49
1:E:101:ARG:HD2	1:E:286:MET:O	2.13	0.49
1:B:33:LEU:O	1:B:36:ILE:HG12	2.14	0.48
1:E:30:LEU:O	1:E:34:MET:HG2	2.14	0.48
1:E:188:LYS:HD3	1:E:188:LYS:HA	1.71	0.47
1:E:71:ARG:NH2	1:E:253:ASP:OD1	2.47	0.47
1:C:91:ILE:O	1:C:95:THR:HG23	2.14	0.47
1:B:42:TYR:HA	1:B:45:TYR:CD1	2.50	0.47
1:E:94:ARG:HH21	1:E:282:ASN:ND2	2.11	0.47
1:D:91:ILE:O	1:D:95:THR:HG23	2.15	0.47
1:B:41:SER:HA	1:B:44:TRP:HD1	1.79	0.46
1:E:130:ARG:HH22	1:E:272:LEU:HB2	1.80	0.46
1:E:71:ARG:HD3	1:E:260:GLU:OE2	2.15	0.46
1:C:191:GLU:OE2	1:D:194:HIS:CE1	2.67	0.46
1:E:75:SER:OG	1:E:260:GLU:HG3	2.16	0.46
1:E:261:ASP:OD2	1:E:264:GLY:HA3	2.16	0.46
1:A:204:THR:O	1:A:206:PRO:HD3	2.16	0.46
1:C:268:ASN:HD21	1:D:208:PRO:HB3	1.81	0.46
1:A:76:TYR:HE1	1:E:207:VAL:HG21	1.80	0.46
1:D:215:LEU:HA	1:D:215:LEU:HD23	1.84	0.45
1:A:91:ILE:O	1:A:95:THR:HG23	2.16	0.45
1:C:168:ILE:HG22	1:C:182:TYR:CD1	2.51	0.45
1:D:168:ILE:HG22	1:D:182:TYR:CD1	2.52	0.45
1:B:209:PHE:O	1:B:213:LEU:HD13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:LYS:HA	1:B:188:LYS:HD3	1.69	0.45
1:C:237:MET:HE2	1:D:52:LEU:HD21	1.99	0.45
1:E:46:GLU:OE2	1:E:51:HIS:HE1	1.99	0.45
1:E:212:THR:O	1:E:216:GLN:HG3	2.17	0.44
1:A:101:ARG:HD2	1:A:286:MET:O	2.17	0.44
1:B:71:ARG:HG3	1:B:215:LEU:HD12	1.98	0.44
1:C:42:TYR:HA	1:C:45:TYR:CD1	2.53	0.44
1:C:268:ASN:ND2	1:D:208:PRO:HB3	2.32	0.44
1:E:89:VAL:CG2	1:E:195:VAL:HG11	2.48	0.44
1:A:149:GLU:OE2	1:A:163:LEU:HD21	2.18	0.44
1:B:208:PRO:HB3	1:D:268:ASN:ND2	2.22	0.44
1:B:126:LYS:HB3	1:B:126:LYS:HE3	1.90	0.44
1:A:215:LEU:HA	1:A:215:LEU:HD23	1.85	0.43
1:A:168:ILE:HG22	1:A:182:TYR:CD1	2.54	0.43
1:B:190:ASP:OD1	1:D:95:THR:HG21	2.19	0.42
1:A:108:HIS:HA	1:A:111:HIS:CG	2.53	0.42
1:C:268:ASN:HD21	1:D:208:PRO:CB	2.33	0.42
1:E:91:ILE:O	1:E:95:THR:HG23	2.19	0.42
1:B:168:ILE:HG22	1:B:182:TYR:CD1	2.54	0.42
1:A:30:LEU:O	1:A:34:MET:HG2	2.20	0.42
1:B:204:THR:O	1:B:206:PRO:HD3	2.20	0.42
1:D:30:LEU:O	1:D:34:MET:HG2	2.20	0.42
1:D:42:TYR:HA	1:D:45:TYR:CD1	2.54	0.42
1:D:213:LEU:HD11	1:D:217:ARG:NH2	2.35	0.42
1:E:202:LEU:HD12	1:E:202:LEU:HA	1.90	0.41
1:D:204:THR:O	1:D:206:PRO:HD3	2.19	0.41
1:D:28:LEU:HD11	1:D:247:TYR:HD1	1.85	0.41
1:A:108:HIS:HA	1:A:111:HIS:ND1	2.36	0.41
1:B:24:ILE:HG22	1:B:28:LEU:HG	2.02	0.41
1:A:285:ASP:O	1:E:159:ARG:NH2	2.53	0.41
1:E:210:ALA:HA	1:E:212:THR:HG22	2.01	0.41
1:A:230:LEU:HB3	1:A:234:LEU:HD12	2.03	0.41
1:B:51:HIS:HD2	1:D:235:HIS:HB2	1.85	0.41
1:A:58:SER:HA	1:E:59:LEU:HD13	2.02	0.41
1:D:271:PRO:HD2	1:D:275:MET:HE3	2.03	0.41
1:A:71:ARG:HG3	1:A:215:LEU:HD12	2.04	0.40
1:A:161:LEU:HA	1:A:161:LEU:HD23	1.90	0.40
1:B:161:LEU:HD23	1:B:161:LEU:HA	1.93	0.40
1:B:85:LEU:HD22	1:B:195:VAL:HA	2.04	0.40

There are no symmetry-related clashes.

## 4.3 Torsion angles

### 4.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	266/301 (88%)	262 (98%)	4 (2%)	0	100	100
1	B	268/301 (89%)	265 (99%)	3 (1%)	0	100	100
1	C	267/301 (89%)	262 (98%)	4 (2%)	1 (0%)	38	72
1	D	266/301 (88%)	263 (99%)	3 (1%)	0	100	100
1	E	268/301 (89%)	263 (98%)	5 (2%)	0	100	100
All	All	1335/1505 (89%)	1315 (98%)	19 (1%)	1 (0%)	55	85

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	267	ALA

### 4.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	229/264 (87%)	228 (100%)	1 (0%)	93	98
1	B	234/264 (89%)	233 (100%)	1 (0%)	93	98
1	C	226/264 (86%)	223 (99%)	3 (1%)	73	93
1	D	228/264 (86%)	227 (100%)	1 (0%)	93	98
1	E	233/264 (88%)	227 (97%)	6 (3%)	51	83
All	All	1150/1320 (87%)	1138 (99%)	12 (1%)	80	95

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	LYS
1	B	71	ARG
1	C	84	ASN
1	C	126	LYS
1	C	190	ASP
1	D	126	LYS
1	E	202	LEU
1	E	207	VAL
1	E	211	TYR
1	E	272	LEU
1	E	285	ASP
1	E	292	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	ASN
1	A	235	HIS
1	A	268	ASN
1	B	43	GLN
1	B	51	HIS
1	B	166	ASN
1	C	51	HIS
1	C	166	ASN
1	C	235	HIS
1	C	268	ASN
1	C	290	HIS
1	D	166	ASN
1	D	268	ASN
1	E	51	HIS
1	E	166	ASN
1	E	235	HIS
1	E	282	ASN

#### 4.3.3 RNA ⓘ

There are no RNA molecules in this entry.

#### 4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

#### 4.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

#### 4.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 15 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

#### 4.7 Other polymers [i](#)

There are no such residues in this entry.

#### 4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 5 Fit of model and data

### 5.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	268/301 (89%)	0.17	18 (6%) 19 14	49, 81, 145, 225	0
1	B	270/301 (89%)	0.35	26 (9%) 9 6	51, 89, 149, 253	0
1	C	269/301 (89%)	0.27	19 (7%) 17 12	52, 86, 164, 217	0
1	D	268/301 (89%)	0.40	22 (8%) 12 9	54, 90, 144, 217	0
1	E	270/301 (89%)	0.31	18 (6%) 19 14	48, 80, 161, 301	0
All	All	1345/1505 (89%)	0.30	103 (7%) 14 10	48, 85, 154, 301	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	265	THR	10.6
1	E	266	ALA	9.6
1	E	47	GLN	9.2
1	E	48	LEU	8.3
1	A	47	GLN	7.4
1	B	220	TYR	7.2
1	C	207	VAL	7.0
1	C	208	PRO	6.8
1	B	291	PRO	6.6
1	E	264	GLY	5.5
1	C	47	GLN	5.3
1	E	269	ASP	5.1
1	A	289	GLN	4.8
1	B	289	GLN	4.8
1	A	175	GLY	4.7
1	D	236	TYR	4.5
1	D	42	TYR	4.4
1	D	265	THR	4.4
1	D	154	SER	4.3
1	B	152	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	211	TYR	4.2
1	A	145	GLU	4.1
1	C	44	TRP	4.0
1	A	48	LEU	3.9
1	B	148	THR	3.8
1	D	44	TRP	3.7
1	D	209	PHE	3.7
1	B	287	THR	3.6
1	B	26	PHE	3.6
1	B	139	ARG	3.5
1	E	292	LEU	3.3
1	A	204	THR	3.3
1	E	26	PHE	3.3
1	B	216	GLN	3.3
1	D	130	ARG	3.2
1	D	204	THR	3.1
1	E	293	PRO	3.1
1	D	264	GLY	3.1
1	B	25	ILE	3.1
1	B	290	HIS	3.0
1	C	24	ILE	3.0
1	A	143	PRO	3.0
1	B	48	LEU	3.0
1	E	204	THR	3.0
1	E	207	VAL	3.0
1	C	30	LEU	2.9
1	E	211	TYR	2.9
1	C	29	LEU	2.9
1	D	266	ALA	2.8
1	E	25	ILE	2.8
1	A	26	PHE	2.8
1	A	63	ALA	2.8
1	A	25	ILE	2.8
1	E	268	ASN	2.7
1	B	149	GLU	2.7
1	A	58	SER	2.7
1	D	43	GLN	2.7
1	D	39	ILE	2.7
1	B	286	MET	2.7
1	A	44	TRP	2.6
1	B	24	ILE	2.6
1	B	138	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	29	LEU	2.6
1	A	209	PHE	2.6
1	D	131	LYS	2.5
1	C	61	GLY	2.5
1	B	146	ARG	2.5
1	B	150	ILE	2.5
1	C	65	ALA	2.5
1	C	266	ALA	2.5
1	D	235	HIS	2.5
1	D	132	THR	2.5
1	E	270	LEU	2.5
1	B	210	ALA	2.4
1	E	44	TRP	2.4
1	C	49	GLY	2.4
1	C	51	HIS	2.4
1	A	174	ALA	2.4
1	C	206	PRO	2.3
1	B	23	LYS	2.3
1	D	289	GLN	2.3
1	C	145	GLU	2.3
1	B	217	ARG	2.3
1	D	173	GLU	2.2
1	D	231	VAL	2.2
1	E	263	PHE	2.2
1	D	129	LEU	2.2
1	A	236	TYR	2.2
1	E	176	LYS	2.1
1	B	153	SER	2.1
1	C	48	LEU	2.1
1	A	55	ALA	2.1
1	B	61	GLY	2.1
1	B	27	ARG	2.1
1	C	43	GLN	2.1
1	A	40	ILE	2.1
1	D	40	ILE	2.1
1	B	36	ILE	2.1
1	C	27	ARG	2.1
1	D	25	ILE	2.1
1	D	272	LEU	2.0
1	C	45	TYR	2.0
1	C	148	THR	2.0



## 5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	304	1/1	0.54	0.09	-0.87	168,168,168,168	0
2	ZN	B	303	1/1	0.83	0.12	-1.09	101,101,101,101	0
2	ZN	C	302	1/1	0.94	0.11	-2.13	107,107,107,107	0
2	ZN	A	302	1/1	0.66	0.21	-	96,96,96,96	0
2	ZN	B	301	1/1	0.56	0.21	-	140,140,140,140	0
2	ZN	E	301	1/1	0.96	0.12	-	88,88,88,88	0
2	ZN	A	303	1/1	0.88	0.10	-	102,102,102,102	0
2	ZN	C	301	1/1	0.74	0.18	-	88,88,88,88	0
2	ZN	D	303	1/1	0.53	0.12	-	189,189,189,189	0
2	ZN	A	301	1/1	0.80	0.15	-	84,84,84,84	0
2	ZN	D	302	1/1	0.85	0.06	-	129,129,129,129	0
2	ZN	C	303	1/1	0.85	0.18	-	97,97,97,97	0
2	ZN	E	302	1/1	0.42	0.18	-	227,227,227,227	0
2	ZN	B	302	1/1	0.71	0.14	-	92,92,92,92	0
2	ZN	D	301	1/1	0.65	0.17	-	117,117,117,117	0

## 5.5 Other polymers [i](#)

There are no such residues in this entry.